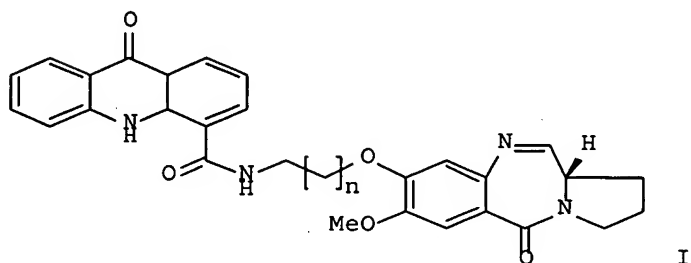
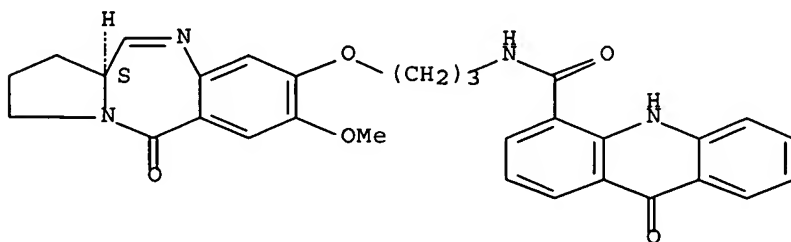


L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN  
 AN 2004:523321 CAPLUS Full-text  
 DN 141:225474  
 TI Synthesis of C8-linked pyrrolo[2,1-c][1,4]benzodiazepine-acridone/acridine hybrids as potential DNA-binding agents  
 AU Kamal, Ahmed; Srinivas, O.; Ramulu, P.; Ramesh, G.; Kumar, P. Praveen  
 CS Division of Organic Chemistry, Indian Institute of Chemical Technology, Hyderabad, 500007, India  
 SO Bioorganic & Medicinal Chemistry Letters (2004), 14(15), 4107-4111  
 CODEN: BMCLE8; ISSN: 0960-894X  
 PB Elsevier Science B.V.  
 DT Journal  
 LA English  
 OS CASREACT 141:225474  
 GI



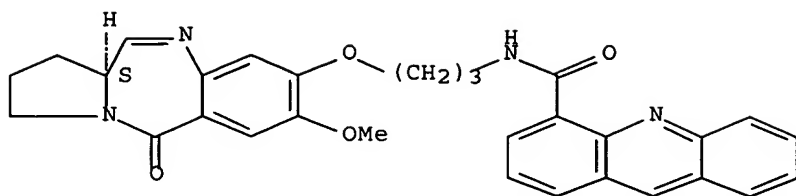
AB Pyrrolobenzodiazepine hybrids linked to acridone/acridine ring systems at C8-position have been designed and prepared. These compounds exhibit significant DNA-binding affinity, and a representative compound (I, n = 2) shows promising in vitro anticancer activity.  
 IT 745014-20-4P 745014-21-5P 745014-22-6P  
 745014-23-7P  
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)  
 (synthesis of C8-linked pyrrolo[2,1-c][1,4]benzodiazepine-acridone/acridine hybrids as potential DNA-binding agents and antitumor agents)  
 RN 745014-20-4 CAPLUS  
 CN 4-Acridinecarboxamide, 9,10-dihydro-9-oxo-N-[3-[[[(11aS)-2,3,5,11a-tetrahydro-7-methoxy-5-oxo-1H-pyrrolo[2,1-c][1,4]benzodiazepin-8-yl]oxy]propyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 745014-21-5 CAPLUS  
 CN 4-Acridinecarboxamide, N-[3-[[[(11aS)-2,3,5,11a-tetrahydro-7-methoxy-5-oxo-1H-pyrrolo[2,1-c][1,4]benzodiazepin-8-yl]oxy]propyl]- (9CI) (CA INDEX NAME)

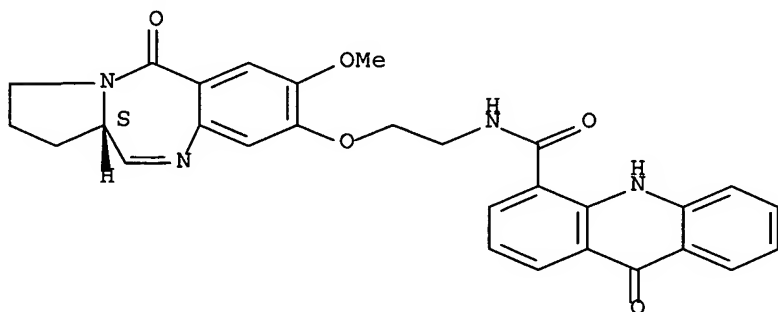
Absolute stereochemistry.



RN 745014-22-6 CAPLUS

CN 4-Acridinecarboxamide, 9,10-dihydro-9-oxo-N-[2-[[[(11aS)-2,3,5,11a-tetrahydro-7-methoxy-5-oxo-1H-pyrrolo[2,1-c][1,4]benzodiazepin-8-yl]oxy]ethyl]- (9CI) (CA INDEX NAME)

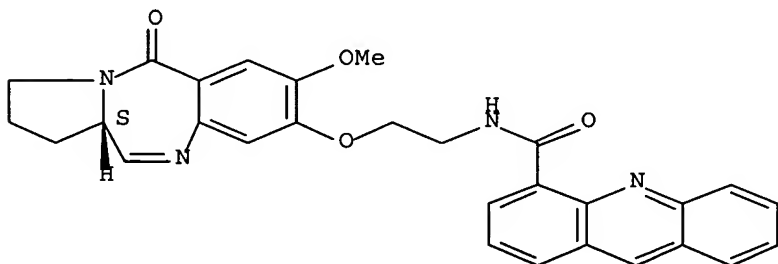
Absolute stereochemistry.



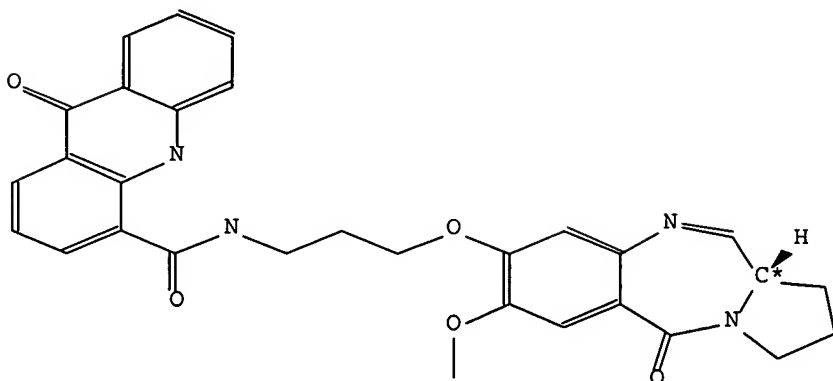
RN 745014-23-7 CAPLUS

CN 4-Acridinecarboxamide, N-[2-[[[(11aS)-2,3,5,11a-tetrahydro-7-methoxy-5-oxo-1H-pyrrolo[2,1-c][1,4]benzodiazepin-8-yl]oxy]ethyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



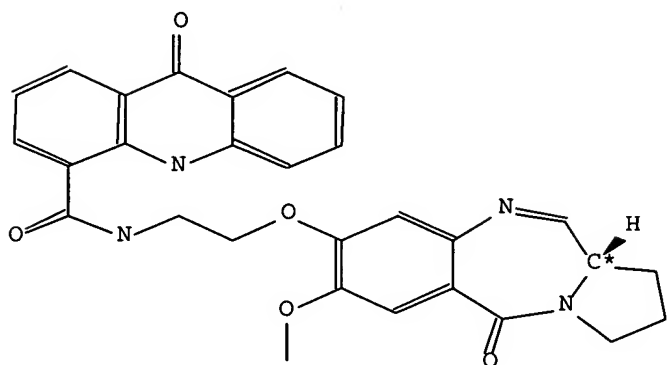
Beilstein Records (BRN):	9824975
Chemical Name (CN):	9-oxo-9,10-dihydro-acridine-4-carboxylic acid <3-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-propyl>-amide
Autonom Name (AUN):	9-oxo-9,10-dihydro-acridine-4-carboxylic acid <3-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-propyl>-amide
Molec. Formula (MF):	C30 H28 N4 O5
Molecular Weight (MW):	524.58
Lawson Number (LN):	29107, 27041, 3131, 289
File Segment (FS):	Stereo compound
Compound Type (CTYPE):	heterocyclic
Constitution ID (CONSID):	8276362
Tautomer ID (TAUTID):	9204410
Entry Date (DED):	2005/01/21
Update Date (DUPD):	2005/01/21



Reference(s):

1. Kamal, Ahmed; Srinivas, O.; Ramulu, P.; Ramesh, G.; Kumar, P. Praveen, Bioorg. Med. Chem. Lett., CODEN: BMCLE8, 14(15), <2004>, 4107 - 4112; BABS-6462175

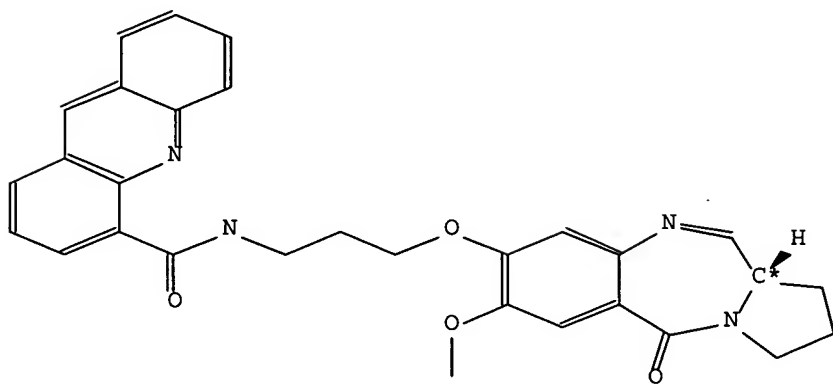
Beilstein Records (BRN):	9824167
Chemical Name (CN):	9-oxo-9,10-dihydro-acridine-4-carboxylic acid <2-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-ethyl>-amide
Autonom Name (AUN):	9-oxo-9,10-dihydro-acridine-4-carboxylic acid <2-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-ethyl>-amide
Molec. Formula (MF):	C29 H26 N4 O5
Molecular Weight (MW):	510.55
Lawson Number (LN):	29107, 27041, 3122, 289
File Segment (FS):	Stereo compound
Compound Type (CTYPE):	heterocyclic
Constitution ID (CONSID):	8275088
Tautomer ID (TAUTID):	9203427
Entry Date (DED):	2005/01/21
Update Date (DUPD):	2005/01/21



Reference(s):

1. Kamal, Ahmed; Srinivas, O.; Ramulu, P.; Ramesh, G.; Kumar, P. Praveen, Bioorg. Med. Chem. Lett., CODEN: BMCLE8, 14(15), <2004>, 4107 - 4112; BABS-6462175

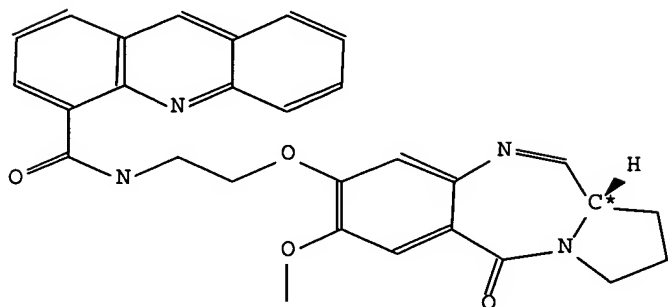
Beilstein Records (BRN): 9823307  
Chemical Name (CN): acridine-4-carboxylic acid  
<3-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-propyl>-amide  
Autonom Name (AUN): acridine-4-carboxylic acid  
<3-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-propyl>-amide  
Molec. Formula (MF): C30 H28 N4 O4  
Molecular Weight (MW): 508.58  
Lawson Number (LN): 29107, 26432, 3131, 289  
File Segment (FS): Stereo compound  
Compound Type (CTYPE): heterocyclic  
Constitution ID (CONSID): 8274346  
Tautomer ID (TAUTID): 9199876  
Entry Date (DED): 2005/01/21  
Update Date (DUPD): 2005/01/21



Reference(s):

1. Kamal, Ahmed; Srinivas, O.; Ramulu, P.; Ramesh, G.; Kumar, P. Praveen, Bioorg. Med. Chem. Lett., CODEN: BMCL E8, 14(15), <2004>, 4107 - 4112; BABS-6462175

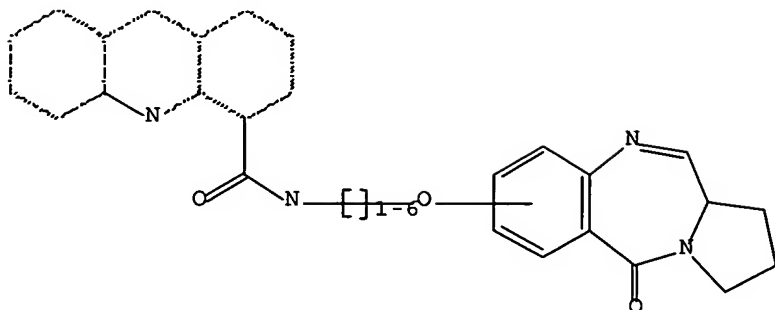
Beilstein Records (BRN): 9821565  
Chemical Name (CN): acridine-4-carboxylic acid  
<2-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-ethyl>-amide  
Autonom Name (AUN): acridine-4-carboxylic acid  
<2-(7-methoxy-5-oxo-2,3,5,11a-tetrahydro-1H-benzo<e>pyrrolo<1,2-a><1,4>diazepin-8-yloxy)-ethyl>-amide  
Molec. Formula (MF): C29 H26 N4 O4  
Molecular Weight (MW): 494.55  
Lawson Number (LN): 29107, 26432, 3122, 289  
File Segment (FS): Stereo compound  
Compound Type (CTYPE): heterocyclic  
Constitution ID (CONSID): 8272807  
Tautomer ID (TAUTID): 9198665  
Entry Date (DED): 2005/01/21  
Update Date (DUPD): 2005/01/21



Reference(s):

1. Kamal, Ahmed; Srinivas, O.; Ramulu, P.; Ramesh, G.; Kumar, P. Praveen, Bioorg. Med. Chem. Lett., CODEN: BMCLE8, 14(15), <2004>, 4107 - 4112; BABS-6462175

=> d l2; d his; log y  
 L2 HAS NO ANSWERS  
 L1 STR



Structure attributes must be viewed using STN Express query preparation.  
 L2 QUE ABB=ON PLU=ON L1

(FILE 'HOME' ENTERED AT 16:55:59 ON 25 MAY 2005)

FILE 'REGISTRY' ENTERED AT 16:56:15 ON 25 MAY 2005

L1 STRUCTURE UPLOADED  
 L2 QUE L1  
 L3 0 S L2  
 L4 4 S L2 FUL

FILE 'CAPLUS' ENTERED AT 16:56:45 ON 25 MAY 2005

L5 1 S L4

FILE 'BEILSTEIN' ENTERED AT 16:57:09 ON 25 MAY 2005

L6 1 S L2  
 L7 4 S L2 FUL  
 L8 4 S L7 NOT L4  
 L9 4 S L7 NOT L5

FILE 'MARPAT' ENTERED AT 16:58:00 ON 25 MAY 2005

L10 0 S L2  
 L11 0 S L2 FUL

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	113.53	400.68
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-0.73

STN INTERNATIONAL LOGOFF AT 16:58:22 ON 25 MAY 2005